Amendments to the claims

 R_1

This listing of claims will replace all prior versions and listings of claims in the application:

1. (Currently amended): <u>A Compounds compound, including stereoisomers</u>, of formula (I) including stereoisomers, prodrugs and pharmaceutically acceptable salts or solvates thereof

$$R_1$$
 N R G G G

or a prodrug, or a pharmaceutically acceptable salt or solvate thereof, wherein the dashed line may represent a double bond;

R is aryl or heteroaryl, each of which may be substituted by 1 to 4 groups J selected from: halogen, C1-C6 alkyl, C1-C6 alkoxy, halo C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, halo C1-C6 alkoxy, -C(O)R₂,

nitro, hydroxy, -NR₃R₄, cyano or a group Z; is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C1-C6 alkoxy, C1-C6 thioalkyl, C2-C6 alkenyl, C2-C6 alkynyl, halo C1-C6

alkyl, halo C1-C6 alkoxy, halogen, NR₃R₄ or cyano;

 R_2 is a C1-C4 alkyl, $-OR_3$ or $-NR_3R_4$;

R₃ is hydrogen or C1-C6 alkyl;

R₄ is hydrogen or C1-C6 alkyl;

is a C1-C6 alkyl, halo C1-C6 alkyl, C1-C6 alkoxy, halo C1-C6 alkoxy, C3-C7 cycloalkyl, hydroxy, halogen, nitro, cyano, -NR₃R₄[[;]], or -C(O)R₂;

R₆ is a C1-C6 alkyl, halo C1-C6 alkyl, C1-C6 alkoxy, halo C1-C6 alkoxy, C3-C7 cycloalkyl, hydroxy, halogen, nitro, cyano, -NR₃R₄[[;]], or -C(O)R₂;

R₇ is hydrogen, C1-C6 alkyl, halogen or halo C1-C6 alkyl;

R₈ is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, NR₃R₄ or cyano;

R₉ is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, NR₃R₄ or cyano;

R₁₀ is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, NR₃R₄ or cyano;

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R₁₁ is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, NR₃R₄ or cyano;

 R_{12} is R_3 or $-C(O)R_2$;

D is CR_8R_9 or is CR_8 when double bonded with G;

G is $CR_{10}R_{11}$ or is CR_{10} when double bonded with D or is CR_{10}

when double bonded with X when X is carbon;

X is carbon or nitrogen;

Y is nitrogen or -CR₇;

W is a 4-8 membered ring, which may be saturated or may

contain one to three double bonds, and

in which:

- one carbon atom is replaced by a carbonyl or S(O)_m; and

- one to four carbon atoms may optionally be replaced by

oxygen, nitrogen or NR_{12} , $S(O)_m$, carbonyl, and such ring may

be further substituted by 1 to 8 R_6 groups;

Z is a 5-6 membered heterocycle, which may be substituted by 1

to 8 R₅ groups or a phenyl ring, which may be substituted by 1

to 4 R₅ groups;

m is an integer from 0 to 2.

2. (Currently amended): <u>A Compounds compound</u> according to claim 1, in which W is selected among from the following groups:

in which:

W1 represents a 1,3-dihydro-2H-imidazol-2-one derivative;

W2 represents a imidazolidin-2-one derivative;

W3 represents a tetrahydropyrimidin-2(1H)-one derivative;

W4 represents a 2,5-dihydro-1,2,5-thiadiazole 1-oxide derivative;

W5 represents a 1,2,5-thiadiazolidine 1-oxide derivative;

W6 represents a 2,5-dihydro-1,2,5-thiadiazole 1,1-dioxide derivative;

W7 represents a 1,2,6-thiadiazinane 1-oxide derivative;

W8 represents a 1,2,6-thiadiazinane 1,1-dioxide derivative;

W9 represents a pyrrolidin-2-one derivative;

W10 represents a 2,5-dihydro-1,2,5-thiadiazolidine 1,1-dioxide derivative;

W11 represents a 1,3-oxazolidin-2-one derivative;

W12 represents a isothiazolidine 1,1-dioxide derivative;

W13 represents a 2(1H)-pyridinone derivative;

W14 represents a 3(2H)-pyridazinone;

W15 represents a 2,3-piperazinedione derivative;

and

q is an integer from 0 to 4[[,]];

n is an integer from 0 to 6[[,]];

p is an integer from 0 to 3; and

m, R₆ and R₁₂ are defined as in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

3. (Currently amended): <u>A Compounds compound according to claim 1, having formula (II)</u>

in which X is nitrogen or carbon and R, R₁, Y, Z, W, D, and G have the meanings as defined in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

- 4. (Currently amended): A Compound according to claim 3, of formula (II), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13, and W14; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- (Currently amended): <u>A Compounds compound</u> according to claim 3 of formula (II), in which Z is selected in the following group from: pyrimidine, pyridine, thiazole, pyrazole, triazole and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 6. (Currently amended): A Compounds compound according to any of claims from 2 to claim 3 of formula (II), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13 and W14 and in which Z is selected from the following heterocyclic groups:

pyrimidine, pyridine, thiazole, pyrazole, triazole and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

7. (Currently amended): A Compound according to any one from claim 1 to claim 6 of formula (IIb), (IIc), (IId), (IIe), (IIf), and or (IIg)

where R, R₁, R₇, Z, W, D, and G have the meanings as defined in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

- 8. (Currently amended): A Compounds compound according to claim 7 of formula (IIb), (IIc), (IId), (IIe), (IIf) and or (IIg), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13 and W14; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 9. (Currently amended): A Compounds compound according to claims claim 7 and 8 of formula (IIb), (IIc), (IId), (IIe), (IIf) and or (IIg), in which Z is selected in the group consisting from: pyrimidine, pyridine, thiazole, pyrazole, triazole and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 10. (Currently amended): A Compounds compound according to any of claims from 7 to 9 claim 7 of formula (IIb), (IIc), (IId), (IIe), (IIf) and or (IIg), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13 and W14 and in which Z is a derivative of the following heterocyclic groups: pyrimidine, pyridine, thiazol, pyrazol, triazol and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 11. (Currently amended): A Compounds compound according to claim 7 of formula (IIr), which corresponds to the compounds a compound of formula (II), where D and G are -CH₂-; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

- 12. (Currently amended): A Compounds compound according to claim 11 of formula (IIr), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13 and W14; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 13. (Currently amended): A Compounds compound according to claims 11 and 12 claim 11 of formula (IIr), in which Z is selected in the group consisting from: pyrimidine, pyridine, thiazol, pyrazol, triazol and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 14. (Currently amended): A Compounds compound according to any of claims from 11 to 13 claim 11 of formula (IIr), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13 and W14 and in which Z is selected in the group consisting from: pyrimidine, pyridine, thiazol, pyrazol, triazol and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 15. (Currently amended): <u>A Compounds compound</u> according to claim 3 of formula (III),

in which Z is a pyrazolyl derivative and R, R₁, R₅, Y, W, D, m and G have the meanings as defined in claim 1 and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

16. (Currently amended): A Compound according to claim 15 of formula (IIIa), (IIIb), (IIIc) and or (IIId),

$$R_{s}$$
 R_{s}
 R_{s}

in which R, R₁, R₅, R₇, R₈, R₉, R₁₀, R₁₁, W, D, m and G have the meanings as defined in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

- 17. (Currently amended): A Compounds compound according to claim 16 of formula (IIIa), (IIIb), (IIIc) and or (IIId), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13, W14 and R, R₁, R₅, R₇, R₈, R₉, R₁₀, R₁₁, and m have the meanings as defined in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 18. (Currently amended): <u>A Compounds compound</u> according to claim 15 of formula (IV),

in which R, R_1 , R_5 , R_6 , R_7 , R_{12} , m, q, D and G have the meanings as defined in claims 1 and 2 and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

19. (Currently amended): A Compounds compound according to claim 18 of formula (IVa), (IVb) and or (IVc),

$$(R_{g})_{q} \qquad (R_{g})_{q} \qquad (R_{g})_{m} \qquad$$

<u>in which</u> R, R₁, R₅, R₆, R₇, R₁₂, m, q, D and G have the meanings as defined in claim 1 and the dashed line may represent a double bond; or a <u>prodrug</u>, or a pharmaceutically acceptable salt or solvate thereof.

20. (Currently amended): A Compounds compound according to claim 3 of formula (V),

<u>in which</u> Z, R, R₁, R₆, q, Y, W, D and G have the meanings as defined in claims 1 and 2, and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

21. (Currently amended): <u>A Compounds compound</u> according to claim 20 of formula (VI),

$$(R_6)_q$$
 N
 R_7
 D
 R_1
 R_1
 R_2
 (VI)

in which Z, R, R₁, R₆, R₇, q, Y, W, D and G have the meanings as defined in claims 1 and 2, and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

22. (Currently amended): A Compounds compound according to claim 21 of formula (VIa), (VIb) and or (VIc),

$$(R_{\theta})_{q}$$

$$Z$$

$$Z$$

$$R_{\theta}$$

$$R$$

$$R_{11}$$

$$R_{10}$$

$$R_{11}$$

$$R_{11}$$

$$R_{11}$$

$$R_{12}$$

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$$R_{12}$$

$$R_{11}$$

$$R_{12}$$

$$R_{12}$$

$$R_{13}$$

$$R_{14}$$

$$R_{15}$$

$$R_{$$

in which R, R_1 , R_6 , R_8 , R_9 , R_{10} , R_{11} , R_{12} , q, D and G have the meanings as defined in claims 1 and 2 and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

- 23. (Currently amended): A Compounds compound according to claim 22 of formula (VIa), (VIb) and or (VIc), in which Z is selected in the group consisting from: pyrimidine, pyridine, thiazol, pyrazol, triazol and phenyl and R, R₁, R₆, R₈, R₉, R₁₀, R₁₁, R₁₂, q, D and G have the meanings as defined in claim 1 and 2; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 24. (Currently amended): A Compounds compound according to any of claims from 1 to 23 of formula (I), (IIb), (IIc), (IId), (IIe), (IIf), (IIg), (IIf), (III), (IIIb), (IIIb), (IIIc), (IIId), (IV), (IVa), (IVb), (IVc), (V), (VI), (VIa), (VIb), (VIe) claim 1, wherein:

R₁ is C1-C3 alkyl group or halo C1-C3 alkyl group, R₇ is hydrogen;

 R_8 , (R_9) , R_{10} , (R_{11}) are hydrogen;

R is an aryl group selected from: 2,4-dichlorophenyl, 2-chloro-4methylphenyl, 2-chloro-4-trifluoromethylphenyl, 2-chloro-4methoxyphenyl, 2,4,5-trimethylphenyl, 2,4-dimethylphenyl, 2-methyl-4methoxyphenyl, 2-methyl-4-ethoxyphenyl, 2-methyl-4-isopropoxyphenyl, 2-methyl-4-hydroxyphenyl, 2-methyl-4-chlorophenyl, 2-methyl-4trifluoromethylphenyl, 2,4-dimethoxyphenyl, 2-methoxy-4trifluoromethylphenyl, 2-methoxy-4-chlorophenyl, 3-methoxy-4chlorophenyl, 2,5-dimethoxy-4-chlorophenyl, 2-methoxy-4isopropylphenyl, 2-methoxy-4-trifluoromethylphenyl, 2-methoxy-4isopropylphenyl, 2-methoxy-4-methylphenyl, 2-trifluoromethyl-4chlorophenyl, 2,4-bis-trifluoromethylphenyl, 2-trifluoromethyl-4methylphenyl, 2-trifluoromethyl-4-methoxyphenyl, 2-difluoromethyl-4methoxyphenyl, 2-bromo-4-isopropylphenyl, 2-methyl-4-cyanophenyl, 2chloro-4-cyanophenyl, 2-trifluoromethyl-4-cyanophenyl, 2trifluoromethoxy-4-cyanophenyl, 2-ethyl-4-cyanophenyl, 2-methyl-4trifluoromethoxyphenyl, 4-methyl-6-dimethylaminopyridin-3-yl, 2,6bismethoxy-pyridin-3-yl, 2-methyl-6-methoxy-pyridin-3-yl, 2trifluoromethyl-6-methoxy-pyridin-3-yl 3-chloro-5-trichloromethyl-

- pyridin-2-yl, 2-methyl-4-(pyrazol-1-yl)-phenyl, 2-methoxy-4-(pyrazol-1-yl)-phenyl, 2,4,6-trimethoxyphenyl, 2-methyl-4,5-benzodioxolyl, and 2-methyl-3,4-benzodioxolyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 25. (Currently amended): A Compounds compound of formula (I), (IIb), (IIe), (IId), (IIe), (IIId), (IIId), (IIId), (IIId), (IIId), (IIId), (IIId), (IV), (IVa), (IVb), (IVc), (IVc), (IVd), (IVd
 - 1-{1-[1-(4-Methoxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}imidazolidin-2-one (compound 1-1);
 - 1-{1-[1-(4-Methoxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}-3-methylimidazolidin-2-one (compound 1-2);
 - 1-{1-[1-(2,4-Dichlorophenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}imidazolidin-2-one (compound 1-3); 1-(1-{1-[2,4-Bis(trifluoromethyl)phenyl]-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-4);
 - 1-{1-[1-(4-Hydroxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}-2-imidazolidinone (compound 1-5);
 - 1-Acetyl-3-(1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-5);
 - 1-Acetyl-3-(1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-6);
 - 1-(1-{1-[4-(Ethyloxy)-2-methylphenyl]-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-7);
 - 1-[1-(6-Methyl-1-{2-methyl-4-[(1-methylethyl)oxy]phenyl}-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl)-1H-pyrazol-3-yl]-2-imidazolidinone (compound 1-8);
 - 1-[1-(6-Methyl-1-{2-methyl-4-[(trifluoromethyl)oxy]phenyl}-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl)-1H-pyrazol-3-yl]-2-imidazolidinone (compound 1-9);
 - 3-Methyl-4-{6-methyl-4-[3-(2-oxo-1-imidazolidinyl)-1H-pyrazol-1-yl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-1-yl}benzonitrile (compound 1-10); 1-(1-{6-Methyl-1-[2-methyl-4-(1H-pyrazol-1-yl)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-11);

- 4-{6-Methyl-4-[3-(2-oxo-1-imidazolidinyl)-1H-pyrazol-1-yl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-1-yl}-3-(trifluoromethyl)benzonitrile (compound 1-12);
- 1-(1-{1-[2-(Difluoromethyl)-4-(methyloxy)phenyl]-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-13);
- 4-{6-Methyl-4-[3-(2-oxo-1-imidazolidinyl)-1H-pyrazol-1-yl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-1-yl}-3-[(trifluoromethyl)oxy]benzonitrile (compound 1-14);
- 3-Ethyl-4-{6-methyl-4-[3-(2-oxo-1-imidazolidinyl)-1H-pyrazol-1-yl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-1-yl}benzonitrile (compound 1-15); 1-(1-{6-Methyl-1-[2-(methyloxy)-4-(1H-pyrazol-1-yl)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound
- 1-16);
- 1-{1-[6-Methyl-1-(6-methyl-1,3-benzodioxol-5-yl)-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}-2-imidazolidinone (compound 1-17);
- 1-(1-{6-Methyl-1-[2,4,6-tris(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-18);
- 1-{1-[6-Methyl-1-(6-methyl-1,3-benzodioxol-5-yl)-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}-2-imidazolidinone (compound 1-19);
- 1-(6-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-2-pyridinyl)-2-imidazolidinone (compound 1-20);
- 1-(4-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-2-pyrimidinyl)-2-imidazolidinone (compound 1-21);
- 1-(2-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-4-pyrimidinyl)-2-imidazolidinone (compound 1-22);
- 1-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-23);
- 1-(1-{2,6-Dimethyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound
- 1-24);
- 1-(3-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}phenyl)-2-imidazolidinone (compound 1-25); 1-(5-Methyl-1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound

1-26);

11-1); and

- 1-[1-(1-{4-[(difluoromethyl)oxy]-2-methylphenyl}-6-methyl-2,3-dihydro-1*H*-pyrrolo[2,3-*b*]pyridin-4-yl)-1*H*-pyrazol-3-yl]-2-imidazolidinone (compound 1-27);
- 1-{1-[1-(4-Methoxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}pyrrolidin-2-one (compound 2-1);
- 1-{1-[1-(4-Methoxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}tetrahydropyrimidin-2(1H)-one (compound 3-1);
- 3-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-1,3-oxazolidin-2-one (compound 4-1);
- Methyl 5-(1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-1,2,5-thiadiazolidine-2-carboxylate 1,1-dioxide) (compound 5-1);
- 4-[3-(1,1-Dioxido-1,2,5-thiadiazolidin-2-yl)-1H-pyrazol-1-yl]-6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridine (compound 5-2).
- 4-[3-(1,1-Dioxido-2-isothiazolidinyl)-1H-pyrazol-1-yl]-6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridine (compound 6-1);
- 3-Methyl-1-(1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2(1H)-pyridinone (compound 7-1);
- 2-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-3(2H)-pyridazinone (compound 8-1);
- 1-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-1,3-dihydro-2H-imidazol-2-one (compound 9-1);
- 1-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 10-1); 1-(6-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-3-pyridinyl)-2-imidazolidinone (compound
- 1-{1-[7-(2,4-Dichlorophenyl)-2-methyl-6,7-dihydro-5H-pyrrolo[2,3-d]pyrimidin-4-yl]-1H-pyrazol-3-yl}-2-pyrrolidinone (compound 11-2); or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 26. (Currently amended): A Process process for the preparation of a compound the compounds of formula (II), starting from a compound compounds of formula (VII), comprising the following steps as in Scheme 1:

Scheme 1

in which

(XXII)

- step a stands for conversion of the leaving group L, selected in a group consisting from: halogen or reactive residue of sulphonic acid (e.g. mesylate, tosylate), preferably chloride, in the compounds (VIII), by reaction with the suitable Z-W derivative;
- step b stands for reduction of the ester group (E) with a suitable reducing agent (such as DIBAl-H) to hydroxy group of compounds (IX);
- step c stands for suitable protection of an NH group eventually present in W group with a P group, such as a p-methoxybenzyl group;
- step d stands for oxidation of the hydroxy group with a suitable oxidizing agent (such as Dess Martin periodinane) to the aldehyde group of compounds (XI);

- steps e + f stands for formation of the aldehyde group of compounds (XIII) by Wittig reaction in the usual conditions, through formation of enol ether followed by acid hydrolysis (step f);
- step g stands for the optional alkylation of the α position of the aldehyde by deprotonation with a suitable base (such as LiN(SiMe₃)₂), followed by the addition of a suitable alkylating agent (such as MeI) to form the alkylated aldehyde of compounds (XIV), (XV);
- step h stands for the conversion of the aldehyde group group by a Grignard reagent (such as MeMgBr) into an alcohol group of compounds (XVI) and (XVIII);
- step i stands for oxidation of the hydroxy group with a suitable oxidizing agent (such as Dess Martin periodinane) to the ketone group of compounds (XVII);
- step j stands for conversion of the hydroxy group in the suitable protecting group of compounds (XIX) (such as TBS: tert-butyldimethylsilyl);
- step k stands for a Buchwald coupling reaction with the suitable amine RNH₂ to give the compounds of formula (XX);
- step 1 stands for the deprotection reaction to give the hydroxy group of compounds (XXI);
- step m stands for intramolecular cyclisation after conversion of the hydroxy group of compounds (XXI) in a suitable leaving group (such as bromide, by reaction with CBr₄ and PPh₃) to give the cyclized compounds (XXII);
- step n stands for the deprotection reaction of the protected NH group eventually present in W group, to give final compounds (II); and
- step o stands for oxidation by a suitable oxidating agent (such as DDQ) in order to give formation of the double bond of compounds (II), when D is CHR₈ and G is CHR₁₀.

27-32. (Cancelled)

- 33. (Currently amended): A pharmaceutical composition comprising a compound according to any of claims from 1 to 25-claim 1,or a prodrug, or a pharmaceutically acceptable salt or solvate thereof, in admixture with one or more physiologically acceptable carriers or excipients.
- 34. (Currently amended): A method for the treatment of a mammal, including man, in particular in the treatment of a condition conditions mediated by CRF (corticotropin-releasing factor), comprising administration of an effective amount of a compound according to any of claims from 1 to 25

claim 1, or a prodrug, or a pharmaceutically acceptable salt or solvate thereof, to a mammal in need thereof.

- 35. (Currently amended): A method, according to claim 34, in the treatment of depression and anxiety, comprising administration of an effective amount of a compound according to any of claims 1 to 25claim 1, or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
- 36. (Currently amended): A method, according to claim 34, in the treatment of IBS (irritable bowel disease) and IBD (inflammatory bowel disease), comprising administration of an effective amount of a compound according to any of claims 1 to 25claim 1, or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.